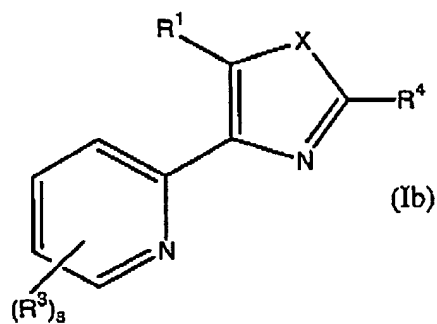
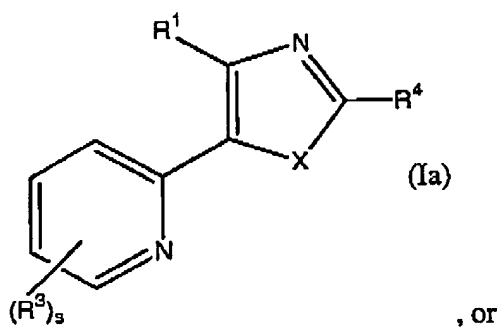


Appl. No. 10/667,187
Amdt. dated 23 February 23, 2006
Reply to Office Action mailed 11/23/05

Amendments to the claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

1. (CURRENTLY AMENDED) A compound of formula (Ia) or (Ib):

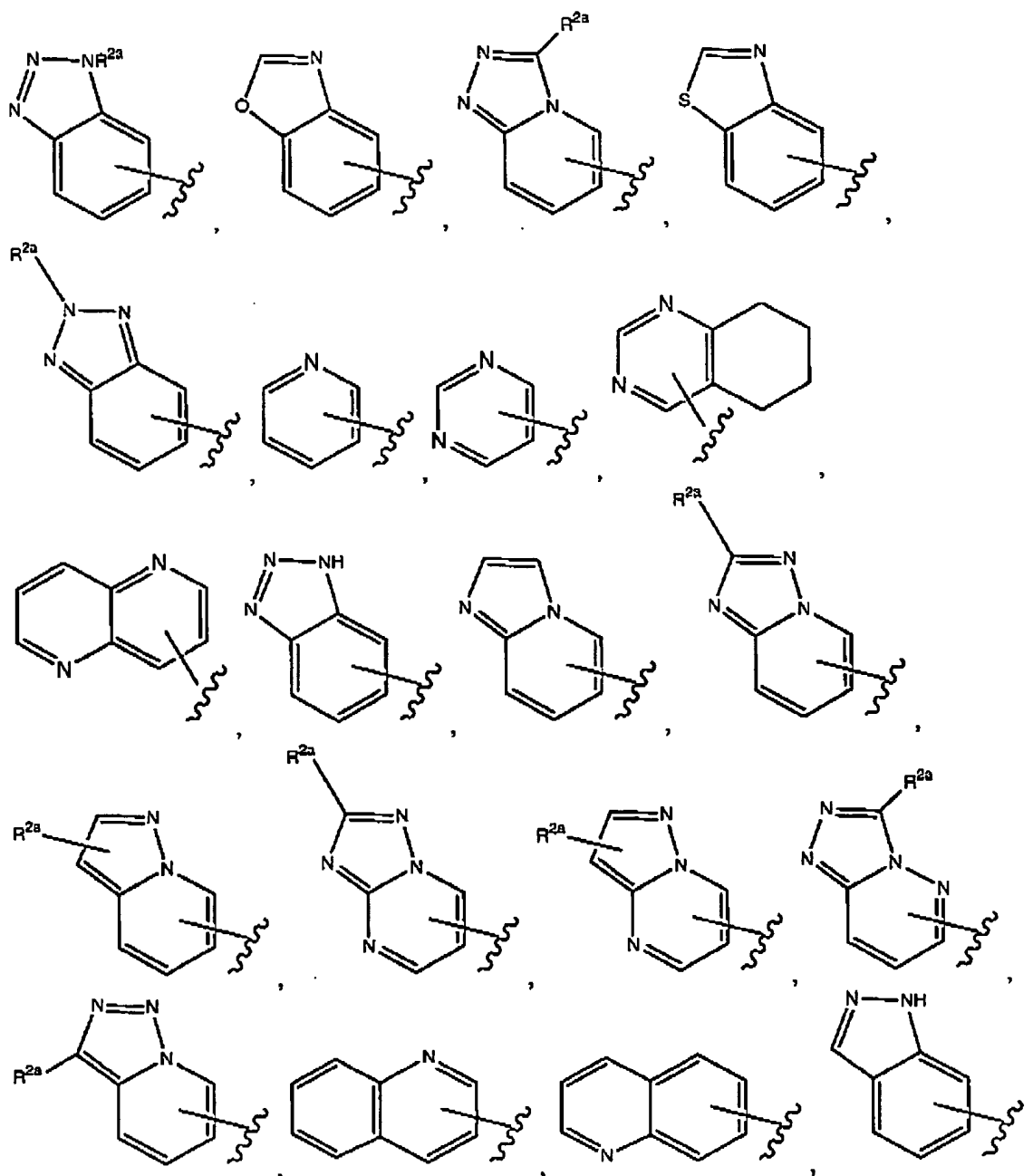


or a pharmaceutically acceptable salt, hydrate, tautomer or solvate thereof, wherein:

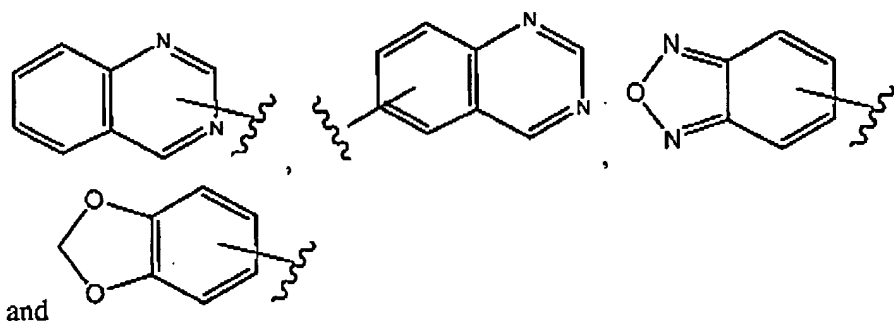
X is O or S;

R¹ is selected from the group consisting of

Appl. No. 10/667,187
Amdt. dated 23 February 23, 2006
Reply to Office Action mailed 11/23/05



Appl. No. 10/667,187
 Amdt. dated 23 February 23, 2006
 Reply to Office Action mailed 11/23/05



where R^{2a} is independently selected from the group consisting of: (C_1-C_6) alkyl, (C_2-C_6) alkenyl, (C_2-C_6) alkynyl, (C_3-C_{10}) cycloalkyl, (C_5-C_{10}) aryl, (C_1-C_6) alkylaryl, amino, carbonyl, carboxyl, (C_2-C_6) acid, (C_1-C_6) ester, (C_5-C_{10}) heteroaryl, (C_5-C_{10}) heterocyclyl, (C_1-C_6) alkoxy, nitro, halo, hydroxyl, (C_1-C_6) alkoxy (C_1-C_6) ester, and where alkyl, alkenyl, alkynyl, cycloalkyl, aryl, amino, (C_2-C_6) acid, (C_1-C_6) ester, heteroaryl, heterocyclyl, and alkoxy of R^{2a} is optionally substituted by at least one moiety independently selected from the group consisting of halo, (C_1-C_6) alkyl, (C_2-C_6) alkenyl, (C_2-C_6) alkynyl, perhalo (C_1-C_6) alkyl, phenyl, (C_3-C_{10}) cycloalkyl, (C_5-C_{10}) heteroaryl, (C_5-C_{10}) heterocyclic, formyl, NC-, (C_1-C_6) alkyl-(C=O)-, phenyl-(C=O)-, HO-(C=O)-, (C_1-C_6) alkyl-O-(C=O)-, (C_1-C_6) alkyl-NH-(C=O)-, $((C_1-C_6)alkyl)_2N-(C=O)-$, phenyl-NH-(C=O)-, phenyl- $[(C_1-C_6)alkyl-N]-(C=O)-$, O_2N- , amino, (C_1-C_6) alkylamino, $((C_1-C_6)alkyl)_2$ -amino, $(C_1-C_6)alkyl-(C=O)-NH-$, $(C_1-C_6)alkyl-(C=O)-[(C_1-C_6)alkyl-N]$, phenyl-(C=O)-NH-, phenyl-(C=O)- $[(C_1-C_6)alkyl-N]$, $H_2N-(C=O)-NH-$, $(C_1-C_6)alkyl-HN-(C=O)-NH-$, $((C_1-C_6)alkyl)_2N-(C=O)-NH-$, $(C_1-C_6)alkyl-HN-(C=O)-[(C_1-C_6)alkyl-N]$, $((C_1-C_6)alkyl)_2N-(C=O)-[(C_1-C_6)alkyl-N]$, phenyl-HN-(C=O)-NH-, $(phenyl)_2N-(C=O)-NH-$, phenyl-HN-(C=O)- $[(C_1-C_6)alkyl-N]$, $(phenyl)_2N-(C=O)-[(C_1-C_6)alkyl-N]$, $(C_1-C_6)alkyl-O-(C=O)-NH-$, $(C_1-C_6)alkyl-O-(C=O)-[(C_1-C_6)alkyl-N]$, phenyl-O-(C=O)-NH-, phenyl-O-(C=O)- $[(alkyl)-N]$, $(C_1-C_6)alkyl-SO_2NH-$, phenyl- SO_2NH- , $(C_1-C_6)alkyl-SO_2-$, phenyl- SO_2- , hydroxy, (C_1-C_6) alkoxy, perhalo (C_1-C_6) alkoxy, phenoxy, $(C_1-C_6)alkyl-(C=O)-O-$, $(C_1-C_6)ester-(C_1-C_6)alkyl-O-$, phenyl-(C=O)-O-, $H_2N-(C=O)-O-$, $(C_1-C_6)alkyl-HN-(C=O)-O-$, $((C_1-C_6)alkyl)_2N-(C=O)-O-$, phenyl-HN-(C=O)-O-, and $(phenyl)_2N-(C=O)-O-$;

Appl. No. 10/667,187
 Amdt. dated 23 February 23, 2006
 Reply to Office Action mailed 11/23/05

wherein R¹ can optionally be further independently substituted with at least one moiety independently selected from the group consisting of: carbonyl, halo, halo(C₁-C₆)alkyl, perhalo(C₁-C₆)alkyl, perhalo(C₁-C₆)alkoxy, (C₁-C₆)alkyl, (C₂-C₆)alkenyl, (C₂-C₆)alkynyl, hydroxy, oxo, mercapto, (C₁-C₆)alkylthio, (C₁-C₆)alkoxy, (C₅-C₁₀)aryl or (C₅-C₁₀)heteroaryl, (C₅-C₁₀)aryloxy or (C₅-C₁₀)heteroaryloxy, (C₅-C₁₀)ar(C₁-C₆)alkyl or (C₅-C₁₀)heteroar(C₁-C₆)alkyl, (C₅-C₁₀)ar(C₁-C₆)alkoxy or (C₅-C₁₀)heteroar(C₁-C₆)alkoxy, HO-(C=O)-, ester, amido, ether, amino, amino(C₁-C₆)alkyl, (C₁-C₆)alkylamino(C₁-C₆)alkyl, di(C₁-C₆)alkylamino(C₁-C₆)alkyl, (C₅-C₁₀)heterocyclyl(C₁-C₆)alkyl, (C₁-C₆)alkyl- and di(C₁-C₆)alkylamino, cyano, nitro, carbamoyl, (C₁-C₆)alkylcarbonyl, (C₁-C₆)alkoxycarbonyl, (C₁-C₆)alkylaminocarbonyl, di(C₁-C₆)alkylaminocarbonyl, (C₅-C₁₀)arylcarbonyl, (C₅-C₁₀)aryloxycarbonyl, (C₁-C₆)alkylsulfonyl, and (C₅-C₁₀)arylsulfonyl;

each R³ is independently selected from the group consisting of: hydrogen, halo, halo(C₁-C₆)alkyl, (C₁-C₆)alkyl, (C₂-C₆)alkenyl, (C₂-C₆)alkynyl, perhalo(C₁-C₆)alkyl, phenyl, (C₅-C₁₀)heteroaryl, (C₅-C₁₀)heterocyclic, (C₃-C₁₀)cycloalkyl, hydroxy, (C₁-C₆)alkoxy, perhalo(C₁-C₆)alkoxy, phenoxy, (C₅-C₁₀)heteroaryl-O-, (C₅-C₁₀)heterocyclic-O-, (C₃-C₁₀)cycloalkyl-O-, (C₁-C₆)alkyl-S-, (C₁-C₆)alkyl-SO₂-, (C₁-C₆)alkyl-NH-SO₂-, O₂N-, NC-, amino, Ph(CH₂)₁₋₆HN-, (C₁-C₆)alkyl HN-, (C₁-C₆)alkylamino, [(C₁-C₆)alkyl]₂-amino, (C₁-C₆)alkyl-SO₂-NH-, amino(C=O)-, aminoO₂S-, (C₁-C₆)alkyl-(C=O)-NH-, (C₁-C₆)alkyl-(C=O)-[((C₁-C₆)alkyl)-N]-, phenyl-(C=O)-NH-, phenyl-(C=O)-[((C₁-C₆)alkyl)-N]-, (C₁-C₆)alkyl-(C=O)-, phenyl-(C=O)-, (C₅-C₁₀)heteroaryl-(C=O)-, (C₅-C₁₀)heterocyclic-(C=O)-, (C₃-C₁₀)cycloalkyl-(C=O)-, HO-(C=O)-, (C₁-C₆)alkyl-O-(C=O)-, H₂N(C=O)-, (C₁-C₆)alkyl-NH-(C=O)-, [(C₁-C₆)alkyl]₂-N-(C=O)-, phenyl-NH-(C=O)-, phenyl-[((C₁-C₆)alkyl)-N]-(C=O)-, (C₅-C₁₀)heteroaryl-NH-(C=O)-, (C₅-C₁₀)heterocyclic-NH-(C=O)-, (C₃-C₁₀)cycloalkyl-NH-(C=O)- and (C₁-C₆)alkyl-(C=O)-O-,

where alkyl, alkenyl, alkynyl, phenyl, heteroaryl, heterocyclic, cycloalkyl, alkoxy, phenoxy, amino of R³ is optionally substituted by at least one substituent independently selected from (C₁-C₆)alkyl, (C₁-C₆)alkoxy, halo(C₁-C₆)alkyl, halo, H₂N-, Ph(CH₂)₁₋₆HN-, and (C₁-C₆)alkylHN-;

Appl. No. 10/667,187
 Amdt. dated 23 February 23, 2006
 Reply to Office Action mailed 11/23/05

s is an integer from one to five;

R^4 is independently selected from the group consisting of: hydrogen, halo, halo(C_1 - C_6)alkyl, (C_1 - C_6)alkyl, (C_2 - C_6)alkenyl, (C_2 - C_6)alkynyl, perhalo(C_1 - C_6)alkyl, phenyl, (C_5 - C_{10})heteroaryl, (C_5 - C_{10})heterocyclic, (C_3 - C_{10})cycloalkyl, hydroxy, (C_1 - C_6)alkoxy, perhalo(C_1 - C_6)alkoxy, phenoxy, (C_5 - C_{10})heteroaryl-O-, (C_5 - C_{10})heterocyclic-O-, (C_3 - C_{10})cycloalkyl-O-, (C_1 - C_6)alkyl-S-, (C_1 - C_6)alkyl-SO₂-, (C_1 - C_6)alkyl-NH-SO₂-, O₂N-, NC-, amino, Ph(CH₂)₁₋₆HN-, (C_1 - C_6)alkylHN-, (C_1 - C_6)alkylamino, [(C_1 - C_6)alkyl]₂-amino, (C_1 - C_6)alkyl-SO₂-NH-, amino(C=O)-, aminoO₂S-, (C_1 - C_6)alkyl-(C=O)-NH-, (C_1 - C_6)alkyl-(C=O)-((C_1 - C_6)alkyl)-N-, phenyl-(C=O)-NH-, phenyl-(C=O)-((C_1 - C_6)alkyl)-N]-, (C_1 - C_6)alkyl-(C=O)-, phenyl-(C=O)-, (C_5 - C_{10})heteroaryl-(C=O)-, (C_5 - C_{10})heterocyclic-(C=O)-, (C_3 - C_{10})cycloalkyl-(C=O)-, HO-(C=O)-, (C_1 - C_6)alkyl-O-(C=O)-, H₂N(C=O)-, (C_1 - C_6)alkyl-NH-(C=O)-, ((C_1 - C_6)alkyl)₂-N-(C=O)-, phenyl-NH-(C=O)-, phenyl-((C_1 - C_6)alkyl)-N]-(C=O)-, (C_5 - C_{10})heteroaryl-NH-(C=O)-, (C_5 - C_{10})heterocyclic-NH-(C=O)-, (C_3 - C_{10})cycloalkyl-NH-(C=O)- and (C_1 - C_6)alkyl-(C=O)-O-,

where alkyl, alkenyl, alkynyl, phenyl, heteroaryl, heterocyclic, cycloalkyl, alkoxy, phenoxy, amino of R^4 is optionally substituted by at least one substituent independently selected from the group consisting of (C_1 - C_6)alkyl, (C_1 - C_6)alkoxy, halo(C_1 - C_6)alkyl, halo, H₂N-, Ph(CH₂)₁₋₆HN-, (C_1 - C_6)alkylHN-, (C_5 - C_{10})heteroaryl and (C_5 - C_{10})heterocyclic;

with the proviso that when R^4 is a substituted phenyl moiety, then (a) R^1 is not naphthyl, phenyl or anthracenyl and (b) if R^1 is a phenyl fused with an aromatic or non-aromatic cyclic ring of 5-7 members wherein said cyclic ring optionally contains up to three heteroatoms independently selected from N, O and S, then the fused cyclic ring of said R^1 moiety is substituted;

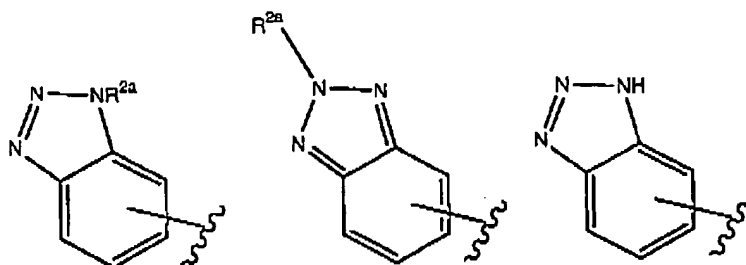
with the proviso that when R^4 is NH₂ and X is S, then R^1 is not an amino-substituted pyridyl or pyrimidinyl moiety;;

with the proviso that when in formula (Ia) R^4 is NH₂ and X is S, then R^1 is not a pyridyl, pyrimidinyl, a naphthyridinyl moiety, or a quinoline moiety that is bonded to the thiazol moiety through the phenyl ring; and

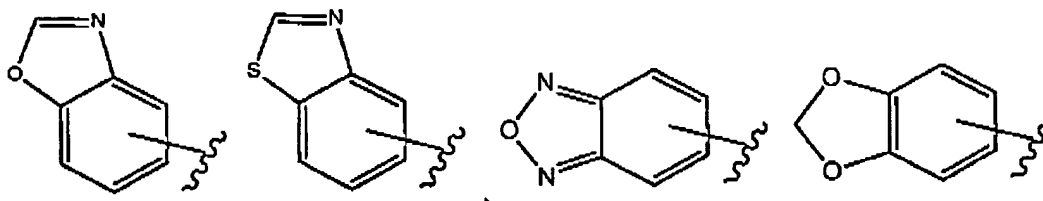
with the proviso that when in formula (Ia) R^4 is CH₃ and X is S, R^1 is not a 3, 4-dimethoxy substituted phenyl moiety.

Appl. No. 10/667,187
 Amdt. dated 23 February 23, 2006
 Reply to Office Action mailed 11/23/05

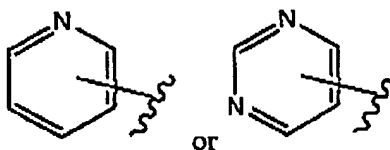
2. (ORIGINAL) A compound of claim 1, wherein R^1 is



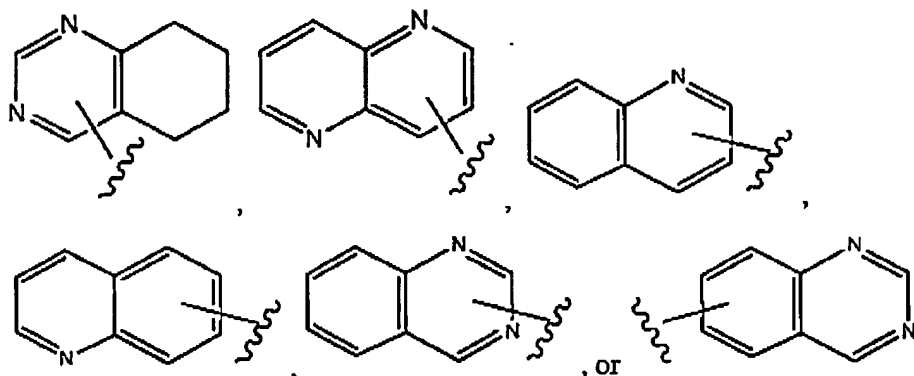
3. (ORIGINAL) A compound of claim 1, wherein R^1 is



4. (ORIGINAL) A compound of claim 1, wherein R^1 is

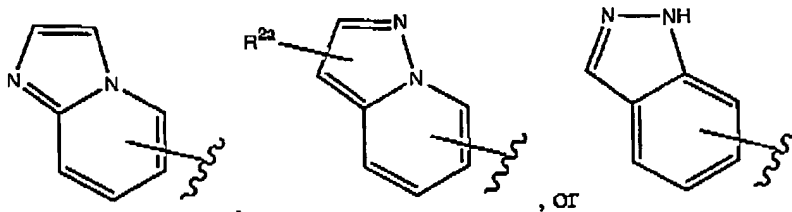


5. (ORIGINAL) A compound of claim 1, wherein R^1 is

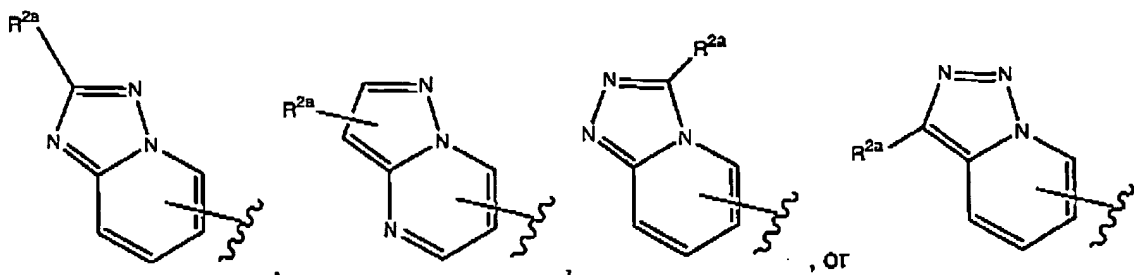


Appl. No. 10/667,187
 Amdt. dated 23 February 23, 2006
 Reply to Office Action mailed 11/23/05

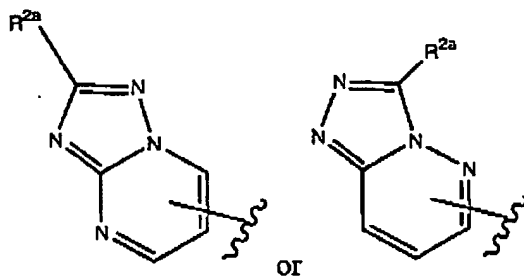
6. (ORIGINAL) A compound of claim 1, wherein R^1 is



7. (ORIGINAL) A compound of claim 1, wherein R^1 is



8. (ORIGINAL) A compound of claim 1, wherein R^1 is



9. (Withdrawn:Non-elected claim) A compound of claim 1, wherein X is O; s is one to two; R^3 is hydrogen or (C_1-C_6) alkyl; and R^4 is H, (C_1-C_6) alkyl, or amino.

10. (ORIGINAL) A compound of claim 1, wherein X is S; s is one to two; R^3 is hydrogen or (C_1-C_6) alkyl; and R^4 is H, (C_1-C_6) alkyl, or amino.

Appl. No. 10/667,187
Amdt. dated 23 February 23, 2006
Reply to Office Action mailed 11/23/05

11. (Previously AMENDED) A pharmaceutical composition comprising a therapeutically effective amount of a compound of claim 1 and a pharmaceutically acceptable carrier.
12. (Previously Presented) A method of treating a TGF-related disease state in an animal or human comprising the step of administering a therapeutically effective amount of a compound of claim 1 to the animal or human suffering from the TGF-related disease state selected from the group consisting of glomerulonephritis, diabetic nephropathy, hepatic fibrosis, pulmonary fibrosis, intimal hyperplasia and restenosis, scleroderma, and dermal scarring.
13. (CANCELED)
14. ((Withdrawn:Non-elected claim) A compound selected from the groups consisting of ;
- 2-(5-Benzo[1,3]dioxol-5-yl-oxazol-4-yl)-6-methyl-pyridine;
 - 2-(5-Benzo[1,3]dioxol-5-yl-oxazol-4-yl)-pyridine;
 - 2-(5-Benzo[1,3]dioxol-5-yl-oxazol-4-yl)-6-methoxy-pyridine;
 - 2-(5-Benzo[1,3]dioxol-5-yl-oxazol-4-yl)-6-trifluoromethyl-pyridine;
 - 2-Methyl-5-[4-(6-methyl-pyridin-2-yl)-oxazol-5-yl]-2H-benzotriazole;
 - 4-[4-(6-Methyl-pyridin-2-yl)-oxazol-5-yl]-quinoline;
 - 1-Methyl-6-[4-(6-methyl-pyridin-2-yl)-oxazol-5-yl]-1H-benzotriazole;
 - 6-(4-Pyridin-2-yl-oxazol-5-yl)-quinoxaline;
 - 6-[4-(6-Methyl-pyridin-2-yl)-oxazol-5-yl]-quinoxaline;
 - 6-[4-(6-Methyl-pyridin-2-yl)-oxazol-5-yl]-quinoline;
 - 6-(4-pyridin-2-yl-oxazol-5-yl)-quinoline;
 - 2-(5-Benzo[1,3]dioxol-5-yl-oxazol-4-yl)-6-ethyl-pyridine;
 - 2-(5-Benzo[1,3]dioxol-5-yl-oxazol-4-yl)-6-propyl-pyridine;
 - 6-[4-(6-Methyl-pyridin-2-yl)-oxazol-5-yl]-benzothiazole;
 - 2-(4-Benzo[1,3]dioxol-5-yl-oxazol-5-yl)-6-methyl-pyridine;
 - 4-[5-(6-Methyl-pyridin-2-yl)-oxazol-4-yl]-quinoline;
 - 1-Methyl-6-[5-(6-methyl-pyridin-2-yl)-oxazol-4-yl]-1H-benzotriazole;
 - 2-Methyl-5-[5-(6-methyl-pyridin-2-yl)-oxazol-4-yl]-2H-benzotriazole;

Appl. No. 10/667,187
Amdt. dated 23 February 23, 2006
Reply to Office Action mailed 11/23/05

6-[5-(6-Methyl-pyridin-2-yl)-oxazol-4-yl]-quinoline;
6-[5-(6-Methyl-pyridin-2-yl)-oxazol-4-yl]-quinoxaline;
2-[5-(6-Methyl-pyridin-2-yl)-oxazol-4-yl]-[1,5]naphthyridine;
{4-[5-(6-Methyl-pyridin-2-yl)-oxazol-4-yl]-pyridin-2-yl}-phenyl-amine;
2-(4-Benzo[1,3]dioxol-5-yl-2-methyl-oxazol-5-yl)-6-methyl-pyridine;
1-Methyl-6-[2-methyl-5-(6-methyl-pyridin-2-yl)-oxazol-4-yl]-1H-benzotriazole;
2-Methyl-5-[2-methyl-5-(6-methyl-pyridin-2-yl)-oxazol-4-yl]-2H-benzotriazole;
6-[2-Methyl-5-(6-methyl-pyridin-2-yl)-oxazol-4-yl]-quinoline;
6-[2-Methyl-5-(6-methyl-pyridin-2-yl)-oxazol-4-yl]-quinoxaline;
2-[2-Methyl-5-(6-methyl-pyridin-2-yl)-oxazol-4-yl]-[1,5]naphthyridine;
{4-[2-Methyl-5-(6-methyl-pyridin-2-yl)-oxazol-4-yl]-pyridin-2-yl}-phenyl-amine;
4-[2-Methyl-5-(6-methyl-pyridin-2-yl)-oxazol-4-yl]-quinoline;
4-Benzo[1,3]dioxol-5-yl-5-(6-methyl-pyridin-2-yl)-thiazol-2-ylamine;
4-(3-Methyl-3H-benzotriazol-5-yl)-5-(6-methyl-pyridin-2-yl)-thiazol-2-ylamine;
4-(2-Methyl-2H-benzotriazol-5-yl)-5-(6-methyl-pyridin-2-yl)-thiazol-2-ylamine;
5-(6-Methyl-pyridin-2-yl)-4-quinolin-6-yl-thiazol-2-ylamine;
5-(6-Methyl-pyridin-2-yl)-4-quinoxalin-6-yl-thiazol-2-ylamine;
5-(6-Methyl-pyridin-2-yl)-4-[1,5]naphthyridin-2-yl-thiazol-2-ylamine;
{4-[2-Amino-5-(6-methyl-pyridin-2-yl)-thiazol-4-yl]-pyridin-2-yl}-phenyl-amine;
5-(6-Methyl-pyridin-2-yl)-4-quinolin-4-yl-thiazol-2-ylamine;
4-(6-Methyl-pyridin-2-yl)-5-quinolin-6-yl-thiazol-2-ylamine;
5-(3-Methyl-3H-benzotriazol-5-yl)-4-(6-methyl-pyridin-2-yl)-thiazol-2-ylamine;
5-(2-Methyl-2H-benzotriazol-5-yl)-4-(6-methyl-pyridin-2-yl)-thiazol-2-ylamine;
5-Benzo[1,3]dioxol-5-yl-4-(6-methyl-pyridin-2-yl)-thiazol-2-ylamine;
4-(6-Methyl-pyridin-2-yl)-5-quinoxalin-6-yl-thiazol-2-ylamine;
4-(6-Methyl-pyridin-2-yl)-5-[1,5]naphthyridin-2-yl-thiazol-2-ylamine;
{4-[2-Amino-4-(6-methyl-pyridin-2-yl)-thiazol-5-yl]-pyridin-2-yl}-phenyl-amine;
4-(6-Methyl-pyridin-2-yl)-5-quinolin-4-yl-thiazol-2-ylamine;
6-[2-Methyl-4-(6-methyl-pyridin-2-yl)-oxazol-5-yl]-quinoline;
1-Methyl-6-[2-methyl-4-(6-methyl-pyridin-2-yl)-oxazol-5-yl]-1H-benzotriazole;
2-Methyl-5-[2-methyl-4-(6-methyl-pyridin-2-yl)-oxazol-5-yl]-2H-benzotriazole;
2-(5-Benzo[1,3]dioxol-5-yl-2-methyl-oxazol-4-yl)-6-methyl-pyridine;

Appl. No. 10/667,187
Amdt. dated 23 February 23, 2006
Reply to Office Action mailed 11/23/05

6-[2-Methyl-4-(6-methyl-pyridin-2-yl)-oxazol-5-yl]-quinoxaline;
2-[2-Methyl-4-(6-methyl-pyridin-2-yl)-oxazol-5-yl]-[1,5]naphthyridine;
{4-[2-Methyl-4-(6-methyl-pyridin-2-yl)-oxazol-5-yl]-pyridin-2-yl}-phenyl-amine;
4-[2-Methyl-4-(6-methyl-pyridin-2-yl)-oxazol-5-yl]-quinoline;
1-Methyl-6-[4-(6-methyl-pyridin-2-yl)-thiazol-5-yl]-1H-benzotriazole;
2-Methyl-5-[4-(6-methyl-pyridin-2-yl)-thiazol-5-yl]-2H-benzotriazole;
2-(5-Benzo[1,3]dioxol-5-yl-thiazol-4-yl)-6-methyl-pyridine;
6-[4-(6-Methyl-pyridin-2-yl)-thiazol-5-yl]-quinoxaline;
2-[4-(6-Methyl-pyridin-2-yl)-thiazol-5-yl]-[1,5]naphthyridine;
{4-[4-(6-Methyl-pyridin-2-yl)-thiazol-5-yl]-pyridin-2-yl}-phenyl-amine;
4-[4-(6-Methyl-pyridin-2-yl)-thiazol-5-yl]-quinoline;
6-[4-(6-Methyl-pyridin-2-yl)-thiazol-5-yl]-quinoline;
1-Methyl-6-[5-(6-methyl-pyridin-2-yl)-thiazol-4-yl]-1H-benzotriazole;
2-Methyl-5-[5-(6-methyl-pyridin-2-yl)-thiazol-4-yl]-2H-benzotriazole;
2-(4-Benzo[1,3]dioxol-5-yl-thiazol-5-yl)-6-methyl-pyridine;
6-[5-(6-Methyl-pyridin-2-yl)-thiazol-4-yl]-quinoxaline;
2-[5-(6-Methyl-pyridin-2-yl)-thiazol-4-yl]-[1,5]naphthyridine;
{4-[5-(6-Methyl-pyridin-2-yl)-thiazol-4-yl]-pyridin-2-yl}-phenyl-amine;
4-[5-(6-Methyl-pyridin-2-yl)-thiazol-4-yl]-quinoline;
6-[5-(6-Methyl-pyridin-2-yl)-thiazol-4-yl]-quinoline;
1-Methyl-6-[2-methyl-4-(6-methyl-pyridin-2-yl)-thiazol-5-yl]-1H-benzotriazole;
2-Methyl-5-[2-methyl-4-(6-methyl-pyridin-2-yl)-thiazol-5-yl]-2H-benzotriazole;
2-(5-Benzo[1,3]dioxol-5-yl-2-methyl-thiazol-4-yl)-6-methyl-pyridine;
6-[2-methyl-4-(6-Methyl-pyridin-2-yl)-thiazol-5-yl]-quinoxaline;
2-[2-methyl-4-(6-Methyl-pyridin-2-yl)-thiazol-5-yl]-[1,5]naphthyridine;
{4-[2-methyl-4-(6-Methyl-pyridin-2-yl)-thiazol-5-yl]-pyridin-2-yl}-phenyl-amine;
4-[2-methyl-4-(6-Methyl-pyridin-2-yl)-thiazol-5-yl]-quinoline;
6-[2-methyl-4-(6-Methyl-pyridin-2-yl)-thiazol-5-yl]-quinoline;
1-Methyl-6-[2-methyl-5-(6-methyl-pyridin-2-yl)-thiazol-4-yl]-1H-benzotriazole;
2-Methyl-5-[2-methyl-5-(6-methyl-pyridin-2-yl)-thiazol-4-yl]-2H-benzotriazole;
2-(4-Benzo[1,3]dioxol-5-yl-2-methyl-thiazol-5-yl)-6-methyl-pyridine;

Appl. No. 10/667,187
Amdt. dated 23 February 23, 2006
Reply to Office Action mailed 11/23/05

6-[2-methyl-5-(6-Methyl-pyridin-2-yl)-thiazol-4-yl]-quinoxaline;
2-[2-methyl-5-(6-Methyl-pyridin-2-yl)-thiazol-4-yl]-[1,5]naphthyridine;
{4-[2-methyl-5-(6-Methyl-pyridin-2-yl)-thiazol-4-yl]-pyridin-2-yl}-phenyl-amine;
4-[2-methyl-5-(6-Methyl-pyridin-2-yl)-thiazol-4-yl]-quinoline and
6-[2-methyl-5-(6-Methyl-pyridin-2-yl)-thiazol-4-yl]-quinoline; or a pharmaceutically
acceptable salt thereof.

15. (Withdrawn: nonelected claim) A pharmaceutical composition comprising a therapeutically
effective amount of a compound of claim 14 and a pharmaceutically acceptable carrier.